

# **WEST VIRGINIA LEGISLATURE**

**2022 REGULAR SESSION**

**Introduced**

## **Senate Bill 666**

BY SENATOR WOODRUM

[Introduced February 17, 2022; referred  
to the Committee on Health and Human Resources]

1 A BILL to amend and reenact §60A-2-204 of the Code of West Virginia, 1931, as amended, all  
 2 generally relating to regulation of controlled substances; adding the active chemicals in  
 3 kratom to Schedule I substances; and adding Delta-8 tetrahydrocannabinol to Schedule I.

*Be it enacted by the Legislature of West Virginia:*

## **ARTICLE 2. STANDARDS AND SCHEDULES.**

### **§60A-2-204. Schedule I.**

1 (a) Schedule I shall consist of the drugs and other substances, by whatever official name,  
 2 common or usual name, chemical name, or brand name designated, listed in this section including  
 3 their isomers, esters, ethers, salts and salts of isomers, esters and ethers, whenever the existence  
 4 of such isomers, esters, ethers and salts is possible within the specific chemical designation.

5 (b) Opiates.

6 Acetyl-alpha-methylfentanyl (N-[1-(1-methyl-2-phenethyl) -4-piperidiny]—  
 7 phenylacetamide);

8 Acetylmethadol;

9 Allylprodine;

10 Alphacetylmethadol (except levoalphacetylmethadol also known as levo-alpha-  
 11 acetylmethadol, levomethadyl acetate, or LAAM);

12 Alphameprodine;

13 Alphamethadol;

14 Alpha-methylfentanyl (N-[1-(alpha-methyl-beta-phenyl) ethyl-4-piperidyl] propionanilide;  
 15 1-(1-methyl-2-phenylethyl)-4-(( propanilido) piperidine);

16 Alpha-methylthiofentanyl (N-[1-methyl-2-(2-thienyl) ethyl- 4-piperidiny]—  
 17 phenylpropanamide);

18 Benzethidine;

19 Betacetylmethadol;

20 Beta-hydroxyfentanyl (N-[1-(2-hydroxy-2-phenethyl) -4- piperidiny]-N-

- 21 phenylpropanamide);
- 22 Beta-hydroxy-3-methylfentanyl (other name: N-[1-(2- hydroxy-2-phenethyl)-3-methyl-4-
- 23 piperidiny]-N-phenylpropanamide);
- 24 Betameprodine;
- 25 Betamethadol;
- 26 Betaprodine;
- 27 Clonitazene;
- 28 Dextromoramide;
- 29 Diampromide;
- 30 Diethylthiambutene;
- 31 Difenoxin;
- 32 Dimenoxadol;
- 33 Dimepheptanol;
- 34 Dimethylthiambutene;
- 35 Dioxaphetyl butyrate;
- 36 Dipipanone;
- 37 Ethylmethylthiambutene;
- 38 Etonitazene;
- 39 Etoxidine;
- 40 Furethidine;
- 41 Hydroxypethidine;
- 42 Ketobemidone;
- 43 Levomoramide;
- 44 Levophenacilmorphan;
- 45 3-Methylfentanyl (N-[3-methyl-1-(2-phenylethyl)-4- piperidyl]-N-phenylpropanamide);
- 46 3-methylthiofentanyl (N-[3-methyl-1-(2-thienyl) ethyl-4- piperidiny]—phenylpropanamide);

- 47 Morpheridine;
- 48 MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
- 49 Noracymethadol;
- 50 Norlevorphanol;
- 51 Normethadone;
- 52 Norpipanone;
- 53 Para-fluorofentanyl (N-(4-fluorophenyl)-N-[1-(2-phenethyl)-4-piperidinyl] propanamide);
- 54 PEPAP(1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine);
- 55 Phenadoxone;
- 56 Phenampromide;
- 57 Phenomorphan;
- 58 Phenoperidine;
- 59 Piritramide;
- 60 Proheptazine;
- 61 Properidine;
- 62 Propiram;
- 63 Racemoramide;
- 64 Thiofentanyl (N-phenyl-N-[1-(2-thienyl)ethyl-4-piperidinyl]-propanamide);
- 65 Tilidine;
- 66 Trimeperidine.
- 67 (c) Opium derivatives:
- 68 Acetorphine;
- 69 Acetyldihydrocodeine;
- 70 Benzylmorphine;
- 71 Codeine methylbromide;
- 72 Codeine-N-Oxide;

- 73           Cyprenorphine;
- 74           Desomorphine;
- 75           Dihydromorphine;
- 76           Drotebanol;
- 77           Etorphine (except HCl Salt);
- 78           Heroin;
- 79           Hydromorphanol;
- 80           Methyldesorphine;
- 81           Methyldihydromorphine;
- 82           Morphine methylbromide;
- 83           Morphine methylsulfonate;
- 84           Morphine-N-Oxide;
- 85           Myrophine;
- 86           Nicocodeine;
- 87           Nicomorphine;
- 88           Normorphine;
- 89           Pholcodine;
- 90           Thebacon.
- 91           (d) Hallucinogenic substances.
- 92           Alpha-ethyltryptamine; some trade or other names: etryptamine; Monase; alpha-ethy-1H-
- 93 indole-3-ethanamine; 3-(2- aminobutyl) indole; alpha-ET; and AET;
- 94           4-bromo-2, 5-dimethoxy-amphetamine; some trade or other names: 4-bromo-2,5-
- 95 dimethoxy-alpha-methylphenethylamine; 4-bromo- 2,5-DMA;
- 96           4-Bromo-2,5-dimethoxyphenethylamine; some trade or other names: 2-(4-bromo-2,5-
- 97 dimethoxyphenyl)-1-aminoethane; alpha- desmethyl DOB; 2C-B, Nexus;
- 98           N-(2-Methoxybenzyl)-4-bromo-2, 5-dimethoxyphenethylamine. The substance has the

99 acronym 25B-NBOMe.

100 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25C-NBOMe)

101 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl) ethanamine (25I-NBOMe)

102 2,5-dimethoxyamphetamine; some trade or other names: 2,5-dimethoxy-alpha-

103 methylphenethylamine; 2,5-DMA;

104 2,5-dimethoxy-4-ethylamphet-amine; some trade or other names: DOET;

105 2,5-dimethoxy-4-(n)-propylthiophenethylamine (other name: 2C-T-7);

106 4-methoxyamphetamine; some trade or other names: 4-methoxy-alpha-

107 methylphenethylamine; paramethoxyamphetamine; PMA;

108 5-methoxy-3, 4-methylenedioxy-amphetamine;

109 4-methyl-2,5-dimethoxy-amphetamine; some trade and other names: 4-methyl-2,5-

110 dimethoxy-alpha-methylphenethylamine; "DOM"; and "STP";

111 3,4-methylenedioxy amphetamine;

112 3,4-methylenedioxymethamphetamine (MDMA);

113 3,4-methylenedioxy-N-ethylamphetamine (also known as ( ethyl-alpha-methyl-3,4

114 (methylenedioxy) phenethylamine, N-ethyl MDA, MDE, MDEA);

115 N-hydroxy-3,4-methylenedioxyamphetamine (also known as ( hydroxy-alpha-methyl-3,4

116 (methylenedioxy) phenethylamine, and ( hydroxy MDA);

117 3,4,5-trimethoxy amphetamine;

118 5-methoxy-N,N-dimethyltryptamine (5-MeO-DMT);

119 Alpha-methyltryptamine (other name: AMT);

120 Bufotenine; some trade and other names: 3-(beta-Dimethylaminoethyl)-5-

121 hydroxyindole;3-(2-dimethylaminoethyl) -5-indolol; N, N-dimethylserotonin; 5-hydroxy-N,N-

122 dimethyltryptamine; mappine;

123 Diethyltryptamine; sometrade and other names: N, N-Diethyltryptamine; DET;

124 Dimethyltryptamine; some trade or other names: DMT;

125 5-Methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT);  
126 Ibogaine; some trade and other names: 7-Ethyl-6, 6 Beta, 7, 8, 9, 10, 12, 13-octahydro-2-  
127 methoxy-6, 9-methano-5H- pyrido [1', 2': 1, 2] azepino [5,4-b] indole; Tabernanthe iboga;  
128 Lysergic acid diethylamide;  
129 Marihuana;  
130 Mescaline;  
131 Parahexyl-7374; some trade or other names: 3-Hexyl -1-hydroxy-7, 8, 9, 10-tetrahydro-6,  
132 6, 9-trimethyl-6H-dibenzo [b,d] pyran; Synhexyl;  
133 Peyote; meaning all parts of the plant presently classified botanically as *Lophophora*  
134 *williamsii* Lemaire, whether growing or not, the seeds thereof, any extract from any part of such  
135 plant, and every compound, manufacture, salts, immediate derivative, mixture or preparation of  
136 such plant, its seeds or extracts;  
137 N-ethyl-3-piperidyl benzilate;  
138 N-methyl-3-piperidyl benzilate;  
139 Psilocybin;  
140 Psilocyn;  
141 Tetrahydrocannabinols; synthetic equivalents of the substances contained in the plant, or  
142 in the resinous extractives of *Cannabis*, sp. and/or synthetic substances, immediate derivatives  
143 and their isomers with similar chemical structure and pharmacological activity such as the  
144 following:  
145 delta-1 Cis or trans tetrahydrocannabinol, and their optical isomers;  
146 delta-6 Cis or trans tetrahydrocannabinol, and their optical isomers;  
147 delta-3,4 Cis or trans tetrahydrocannabinol, and its optical isomers;  
148 delta-8 tetrahydrocannabinol and its optical isomers;  
149 (Since nomenclature of these substances is not internationally standardized, compounds  
150 of these structures, regardless of numerical designation of atomic positions covered.)

151 Ethylamine analog of phencyclidine; some trade or other names: N-ethyl-1-  
152 phenylcyclohexylamine, (1-phenylcyclohexyl) ethylamine, N-(1-phenylcyclohexyl) ethylamine,  
153 cyclohexamine, PCE;

154 Pyrrolidine analog of phencyclidine; some trade or other names: 1-(1-phenylcyclohexyl)-  
155 pyrrolidine, PCPy, PHP;

156 Thiophene analog of phencyclidine; some trade or other names: 1-[1-(2-thienyl)-  
157 cyclohexyl]-piperidine, 2-thienyl analog of phencyclidine; TCP, TCP;

158 1[1-(2-thienyl)cyclohexyl]pyrrolidine; some other names: TCPy.

159 4-methylmethcathinone (Mephedrone);

160 3,4-methylenedioxypropylvalerone (MDPV);

161 2-(2,5-Dimethoxy-4-ethylphenyl)ethanamine (2C-E);

162 2-(2,5-Dimethoxy-4-methylphenyl)ethanamine (2C-D)

163 2-(4-Chloro-2,5-dimethoxyphenyl)ethanamine (2C-C)

164 2-(4-Iodo-2,5-dimethoxyphenyl)ethanamine (2C-I)

165 2-[4-(Ethylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-2)

166 2-[4-(Isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4)

167 2-(2,5-Dimethoxyphenyl)ethanamine (2C-H)

168 2-(2,5-Dimethoxy-4-nitro-phenyl)ethanamine (2C-N)

169 2-(2,5-Dimethoxy-4-(n)-propylphenyl)ethanamine (2C-P)

170 3,4-Methylenedioxy-N-methylcathinone (Methylone)

171 2,5-dimethoxy-4-(n)-propylthiophenethylamine (2C-T-7, its optical isomers, salts and  
172 salts of isomers)

173 5-methoxy-N,N-dimethyltryptamine some trade or other names: 5-methoxy-3-[2-  
174 (dimethylamino)ethyl]indole; 5-MeO-DMT(5-MeO-DMT)

175 Alpha-methyltryptamine (other name: AMT)

176 5-methoxy-N,N-diisopropyltryptamine (other name: 5-MeO-DIPT)

177 Synthetic Cannabinoids as follows:

178 2-[(1R,3S)-3-hydroxycyclohexyl]-5-(2-methyloctan-2-yl)phenol {also known as CP  
179 47,497 and homologues};

180 rel-2-[(1S,3R)-3-hydroxycyclohexyl]-5-(2-methylnonan-2-yl)phenol {also known as CP  
181 47,497-C8 homolog};

182 [(6aR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-  
183 tetrahydrobenzo[c]chromen-1-ol] {also known as HU-210};  
184 (dexanabinol);

185 (6aS,10aS)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-  
186 tetrahydrobenzo[c]chromen-1-ol {also known as HU-211};

187 1-Pentyl-3-(1-naphthoyl)indole {also known as JWH-018};

188 1-Butyl-3-(1-naphthoyl)indole {also known as JWH-073};

189 (2-methyl-1-propyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-015};

190 (1-hexyl-1H-indol-3-yl)-1-naphthalenyl-methanone {also known as JWH-019};

191 [1-[2-(4-morpholinyl) ethyl]-1H-indol-3-yl]-1-naphthalenyl-methanone {also known as  
192 JWH-200};

193 1-(1-pentyl-1H-indol-3-yl)-2-(3-hydroxyphenyl)-ethanone {also known as JWH-250};

194 2-[(1S,2S,5S)-5-hydroxy-2-(3-hydroxypropyl)cyclohexyl]-5-(2-methyloctan-2-yl)phenol  
195 {also known as CP 55,940};

196 (4-methyl-1-naphthalenyl)(1-pentyl-1H-indol-3-yl)-methanone {also known as JWH-122};

197 (4-methyl-1-naphthalenyl)(1-pentyl-1H-indol-3-yl)-methanone {also known as JWH-398};

198 (4-methoxyphenyl)(1-pentyl-1H-indol-3-yl)methanone {also known as RCS-4};

199 1-(1-(2-cyclohexylethyl)-1H-indol-3-yl)-2-(2-methoxyphenyl) ethanone {also known as  
200 RCS-8};

201 1-pentyl-3-[1-(4-methoxynaphthoyl)]indole (JWH-081);

202 1-(5-fluoropentyl)-3-(1-naphthoyl)indole (AM2201); and

203 1-(5-fluoropentyl)-3-(2-iodobenzoyl)indole (AM694).

204 Synthetic cannabinoids:

205 CP 47,497 AND homologues, 2-[(1R,3S)-3-Hydroxycyclohexyl]-5-(2-methyloctan-2-

206 YL)phenol);

207 HU-210, [(6AR,10AR)-9-(hydroxymethyl)-6,6-dimethyl-3-(2-Methyloctan-2-YL)-6A,7,10,

208 10A-tetrahydrobenzo[C] chromen-1-OL)];

209 HU-211, (dexanabinol, (6AS,10AS)-9-(hydroxymethyl)-6,6-Dimethyl-3-(2-methyloctan-2-

210 YL)-6A,7,10,10atetrahydrobenzo[ C]chromen-1-OL);

211 7-hydroxymitragynine:

212 JWH-018, 1-pentyl-3-(1-naphthoyl)indole;

213 JWH-019, 1-hexyl-3-(1-naphthoyl)indole;

214 JWH-073, 1-butyl-3-(1-naphthoyl)indole;

215 JWH-200, (1-(2-morpholin-4-ylethyl)indol-3-yl)- Naphthalen-1-ylmethanone;

216 JWH-250, 1-pentyl-3-(2-methoxyphenylacetyl)indole.]

217 Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate (5F-

218 ADB);

219 Methyl 2-(1-(5-fluoropentyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (5F-AMB);

220 Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3-methylbutanoate (FUB-

221 AMB);

222 N-(adamantan-1-yl)-1-(5-fluoropentyl)-1H-indazole-3-carboxamide (5F-APINACA);

223 N-(1-amino-3,3-dimethyl-1-oxobutan-2-yl)-1-(4-fluorobenzyl)-1H-indazole-3-carboxamide

224 (ADB-FUBINACA);

225 Methyl 2-(1-(cyclohexylmethyl)-1H-indole-3-carboxamido)-3,3-dimethylbutanoate

226 (MDMB-CHMICA);

227 Methyl 2-(1-(4-fluorobenzyl)-1H-indazole-3-carboxamido)-3,3-dimethylbutanoate

228 (MDMB-FUBINACA);

- 229 Mitragynine;
- 230 Tetrahydrocannabinols:
- 231 DELTA-1 CIS OR trans tetrahydrocannabinol and their Optical isomers.
- 232 DELTA-6 CIS OR trans tetrahydrocannabinol and their optical isomers.
- 233 DELTA-3,4 CIS or their trans tetrahydrocannabinol and their optical isomers.
- 234 Synthetic Phenethylamines
- 235 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25I-NBOMe/ 2C-I-
- 236 NBOMe);
- 237 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25C-NBOMe/2C-C-
- 238 NBOMe);
- 239 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine (25B-NBOMe/ 2C-B-
- 240 NBOMe);
- 241 Synthetic Opioids (including their isomers, esters, ethers, salts and salts of isomers, esters
- 242 and ethers):
- 243 N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide (acetyl fentanyl);
- 244 furanyl fentanyl;
- 245 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide (also known as U-
- 246 47700);
- 247 N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide, also known as N-(1-
- 248 phenethylpiperidin-4-yl)-N-phenylbutanamide, (butyryl fentanyl);
- 249 N-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-N-phenylpropionamide, also known
- 250 as N-[1-[2-hydroxy-2-(2-thienyl)ethyl]-4-piperidinyll]-N-phenylpropanamide, (beta-
- 251 hydroxythiofentanyl).
- 252 N-(1-phenethylpiperidin-4-yl)-N-phenylacrylamide (acryl fentanyl)
- 253 N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide (isobutyryl fentanyl)
- 254 N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide (cyclopropyl fentanyl)

255 2-(2,4-dichlorophenyl)-N-((1S,2S)-2-(dimethylamino)cyclohexyl)-N-methylacetamide

256 (also known as U-48800)

257 Trans-3,4-dichloro-N-[2-(diethylamino)cyclohexyl]-N-methyl-benzamide (also known as

258 U-49900)

259 Trans-3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methyl-benzeneacetamide (also

260 known as U-51754)

261 Opioid Receptor Agonist

262 AH-7921 (3,4-dichloro-N-(1dimethylamino)cyclohexylmethyl]benzamide).

263 Naphthoylindoles or any compound containing a 3-(1-Naphthoyl) indole structure with

264 substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole

265 ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall

266 include the following:

267 JWH 015;

268 JWH 018;

269 JWH 019;

270 JWH 073;

271 JWH 081;

272 JWH 122;

273 JWH 200;

274 JWH 210;

275 JWH 398;

276 AM 2201;

277 WIN 55,212.

278 Naphylmethylindoles or any compound containing a 1-hindol-3-yl-(1-naphthyl) methane

279 structure with a substitution at the nitrogen atom of the indole ring whether or not further

280 substituted in the indole ring to any extent and whether or not substituted in the naphthyl ring to

281 any extent. This shall include, but not be limited to, JWH 175 and JWH 184.

282 Naphthoylpyrroles or any compound containing a 3-(1- Naphthoyl) pyrrole structure with  
283 substitution at the nitrogen atom of the pyrrole ring whether or not further substituted in the pyrrole  
284 ring to any extent and whether or not substituted in the naphthyl ring to any extent. This shall  
285 include, but not be limited to, JWH 147 and JWH 307.

286 Naphthylmethylindenes or any compound containing a Naphthylideneindene structure  
287 with substitution at the 3- Position of the indene ring whether or not further substituted in the  
288 indene ring to any extent and whether or not substituted in the naphthyl ring to any extent. This  
289 shall include, but not be limited to, JWH 176.

290 Phenylacetylindoles or any compound containing a 3- Phenylacetylindole structure with  
291 substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole  
292 ring to any extent and whether or not substituted in the phenyl ring to any extent. This shall include  
293 the following:

294 RCS-8, SR-18 OR BTM-8;

295 JWH 250;

296 JWH 203;

297 JWH 251;

298 JWH 302.

299 Cyclohexylphenols or any compound containing a 2-(3- hydroxycyclohexyl) phenol  
300 structure with a substitution at the 5-position of the phenolic ring whether or not substituted in the  
301 cyclohexyl ring to any extent. This shall include the following:

302 CP 47,497 and its homologues and analogs;

303 Cannabicyclohexanol;

304 CP 55,940.

305 Benzoylindoles or any compound containing a 3-(benzoyl) indole structure with  
306 substitution at the nitrogen atom of the indole ring whether or not further substituted in the indole

307 ring to any extent and whether or not substituted in the phenyl ring to any extent. This shall include  
308 the following:

309 AM 694;

310 Pravadoline WIN 48,098;

311 RCS 4;

312 AM 679.

313 [2,3-dihydro-5 methyl-3-(4-morpholinylmethyl)pyrrolo [1,2,3-DE]-1, 4-benzoxazin-6-YL]-1-  
314 naphthalenymethanone. This shall include WIN 55,212-2.

315 Dibenzopyrans or any compound containing a 11-hydroxydelta 8-tetrahydrocannabinol  
316 structure with substitution on the 3-pentyl group. This shall include HU-210, HU-211, JWH 051  
317 and JWH 133.

318 Adamantoylindoles or any compound containing a 3-(-1- Adamantoyl) indole structure with  
319 substitution at the nitrogen atom of the indole ring whether or not further substituted in the  
320 adamantoyl ring system to any extent. This shall include AM1248.

321 Tetramethylcyclopropylindoles or any compound containing A 3-  
322 tetramethylcyclopropylindole structure with substitution at the nitrogen atom of the indole ring  
323 whether or not further substituted in the indole ring to any extent and whether or not substituted  
324 in the tetramethylcyclopropyl ring to any extent. This shall include UR-144 and XLR-11.

325 N-(1-Adamantyl)-1-pentyl-1h-indazole-3-carboxamide. This shall include AKB48.

326 Any other synthetic chemical compound that is a Cannabinoid receptor type 1 agonist as  
327 demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV and  
328 V, not federal Food and Drug Administration approved drug or used within legitimate, approved  
329 medical research. Since nomenclature of these substances is not internationally standardized,  
330 any immediate precursor or immediate derivative of these substances shall be covered.

331 Tryptamines:

332 5- methoxy- N- methyl-N-isopropyltryptamine (5-MeO-MIPT)

- 333 4-hydroxy-N,N-diisopropyltryptamine (4-HO-DiPT)
- 334 4-hydroxy-N-methyl-N-isopropyltryptamine (4-HO-MiPT)
- 335 4-hydroxy-N-methyl-N-ethyltryptamine (4-HO-MET)
- 336 4-acetoxy-N,N-diisopropyltryptamine (4-AcO-DiPT)
- 337 5-methoxy- $\alpha$ -methyltryptamine (5-MeO-AMT)
- 338 4-methoxy-N,N-Dimethyltryptamine (4-MeO-DMT)
- 339 4-hydroxy Diethyltryptamine (4-HO-DET)
- 340 5-methoxy-N,N-diallyltryptamine (5-MeO-DALT)
- 341 4-acetoxy-N,N-Dimethyltryptamine (4-AcO DMT)
- 342 4-hydroxy Diethyltryptamine (4-HO-DET)
- 343 (e) Depressants.
- 344 Mecloqualone;
- 345 Methaqualone.
- 346 (f) Stimulants.
- 347 Aminorex; some other names: aminoxaphen; 2-amino-5-phenyl-2-oxazoline; or 4,5-
- 348 dihydro-5-phenyl-2-oxazolamine;
- 349 Cathinone; some trade or other names: 2-amino-1-phenyl-1-propanone, alpha-
- 350 aminopropiophenone, 2-aminopropiophenone and norephedrone;
- 351 Fenethylamine;
- 352 Methcathinone, its immediate precursors and immediate derivatives, its salts, optical
- 353 isomers and salts of optical isomers; some other names: (2-(methylamino)-propiofenone; alpha-
- 354 (methylamino)propiofenone; 2-(methylamino)-1-phenylpropan-1-one; alpha---
- 355 methylaminopropiophenone; monomethylpropion; 3,4-methylenedioxypropiofenone and/or
- 356 mephedrone; 3,4-methylenedioxypropiofenone (MPVD); ephedrone; N-methylcathinone;
- 357 methylcathinone; AL-464; AL-422; AL-463 and UR1432;
- 358 (+-) cis-4-methylaminorex; ((+)-cis-4,5-dihydro-4-methyl-5-phenyl-2-oxazolamine);

- 359 N-ethylamphetamine;
- 360 N,N-dimethylamphetamine; also known as N,N-alpha- trimethyl-benzeneethanamine;
- 361 N,N-alpha-trimethylphenethylamine.
- 362 Alpha-pyrrolidinopentiophenone, also known as alpha-PVP, optical isomers, salts and
- 363 salts of isomers.
- 364 Substituted amphetamines:
- 365 2-Fluoroamphetamine
- 366 3-Fluoroamphetamine
- 367 4-Fluoroamphetamine
- 368 2-chloroamphetamine
- 369 3-chloroamphetamine
- 370 4-chloroamphetamine
- 371 2-Fluoromethamphetamine
- 372 3-Fluoromethamphetamine
- 373 4-Fluoromethamphetamine
- 374 4-chloromethamphetamine
- 375 (g) Temporary listing of substances subject to emergency scheduling. Any material,
- 376 compound, mixture or preparation which contains any quantity of the following substances:
- 377 N-[1-benzyl-4-piperidyl]-N-phenylpropanamide (benzylfentanyl), its optical isomers, salts,
- 378 and salts of isomers.
- 379 N-[1-(2-thienyl)methyl-4-piperidyl]-N-phenylpropanamide (thenylfentanyl), its optical
- 380 isomers, salts and salts of isomers.
- 381 N-benzylpiperazine, also known as BZP.
- 382 Cyclopentyl fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide);
- 383 4-fluorobutyryl fentanyl (N-(4-fluorophenyl)-N-[1-(2-phenylethyl)piperidin-4-yl]-
- 384 butyramide);

385 Isobutyryl fentanyl (2-methyl-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]-propanamide);  
386 Methoxyacetyl fentanyl (2-methoxy-N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]-  
387 acetamide);  
388 3-methylbutyryl fentanyl (N-[3-methyl-1-(2-phenylethyl)piperidin-4-yl]-N-  
389 phenylbutyramide);  
390 4-methoxybutyryl fentanyl (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-  
391 yl)butyramide);  
392 Ocfentanil (N-(2-fluorophenyl)-2-methoxy-N-[1-(2-phenylethyl)piperidin-4-yl]-acetamide);  
393 Tetrahydrofuran fentanyl (N-(1-phenethylpiperidin-4-yl)-N-phenyltetrahydrofuran-2-  
394 carboxamide);  
395 Valeryl fentanyl (N-phenyl-N-[1-(2-phenylethyl)piperidin-4-yl]pentanamide).

396 (h) The following controlled substances are included in Schedule I:

397 Synthetic Cathinones or any compound, except bupropion or compounds listed under a  
398 different schedule, or compounds used within legitimate and approved medical research,  
399 structurally derived from 2- Aminopropan-1-one by substitution at the 1-position with Monocyclic  
400 or fused polycyclic ring systems, whether or not the compound is further modified in any of the  
401 following ways:

402 By substitution in the ring system to any extent with Alkyl, alkylenedioxy, alkoxy, haloalkyl,  
403 hydroxyl or halide Substituents whether or not further substituted in the ring system by one or  
404 more other univalent substituents.

405 By substitution at the 3-position with an acyclic alkyl substituent.

406 By substitution at the 2-amino nitrogen atom with alkyl, dialkyl, benzyl or methoxybenzyl  
407 groups.

408 By inclusion of the 2-amino nitrogen atom in a cyclic structure.

409 Any other synthetic chemical compound that is a Cannabinoid receptor type 1 agonist as  
410 demonstrated by binding studies and functional assays that is not listed in Schedules II, III, IV and

- 411 V, not federal Food and Drug Administration approved drug or used within legitimate, approved
- 412 medical research.

NOTE: The purpose of this bill is to add Kratom and Delta-8 THC to the list of Schedule 1 controlled substances.

Strike-throughs indicate language that would be stricken from a heading or the present law and underscoring indicates new language that would be added.